

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122221\
 Data File : BN018168.D
 Acq On : 23 Dec 2021 01:34
 Operator : CG/JU
 Sample : PB141476BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS476

Quant Time: Dec 23 02:14:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 14:08:04 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	3157	0.400	ng/ul	0.00
4) Naphthalene-d8	10.392	136	10242	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.252	164	6649	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.994	188	14862	0.400	ng/ul	0.00
17) Chrysene-d12	21.190	240	14587	0.400	ng/ul	0.00
23) Perylene-d12	23.414	264	13673	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.075	96	2473	0.758	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.992	152	5662	0.402	ng/ul	0.00
18) Fluoranthene-d10	19.030	212	16721	0.423	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	7364	1.866	ng/ul#	95
5) Naphthalene	10.442	128	11415	0.392	ng/ul	100
7) 2-Methylnaphthalene	12.069	142	7258	0.388	ng/ul	99
8) 1-Methylnaphthalene	12.284	142	7503	0.386	ng/ul	100
10) Acenaphthylene	13.970	152	10285	0.391	ng/ul	99
11) Acenaphthene	14.313	153	8954	0.386	ng/ul	100
12) Fluorene	15.303	166	10406	0.388	ng/ul	99
14) Pentachlorophenol	16.664	266	1588	0.569	ng/ul	98
15) Phenanthrene	17.036	178	18942	0.390	ng/ul	100
16) Anthracene	17.129	178	15100	0.385	ng/ul	99
19) Fluoranthene	19.058	202	23595	0.421	ng/ul	99
20) Pyrene	19.421	202	24614	0.420	ng/ul	97
21) Benzo(a)anthracene	21.172	228	18971	0.391	ng/ul	99
22) Chrysene	21.225	228	25133	0.423	ng/ul	99
24) Benzo(b)fluoranthene	22.744	252	22367	0.427	ng/ul	100
25) Benzo(k)fluoranthene	22.791	252	28041	0.430	ng/ul	100
26) Benzo(a)pyrene	23.317	252	20392	0.411	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.672	276	25516	0.391	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.695	278	19263	0.389	ng/ul	99
29) Benzo(g,h,i)perylene	26.363	276	21148	0.392	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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